

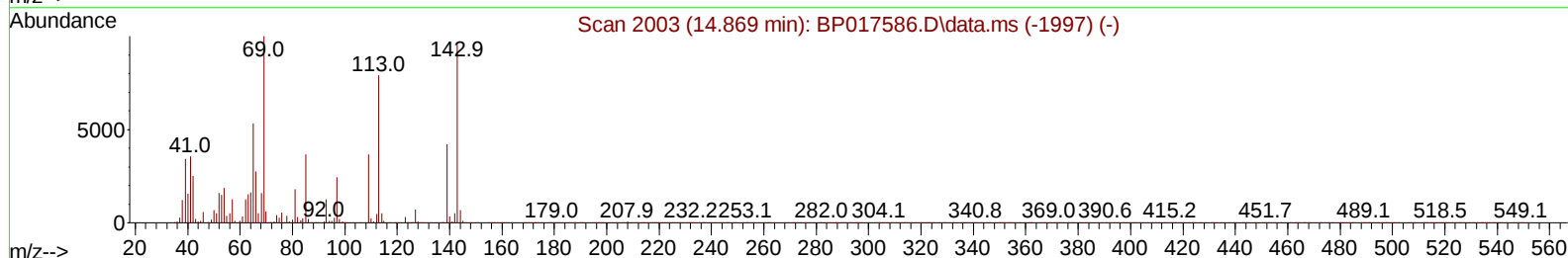
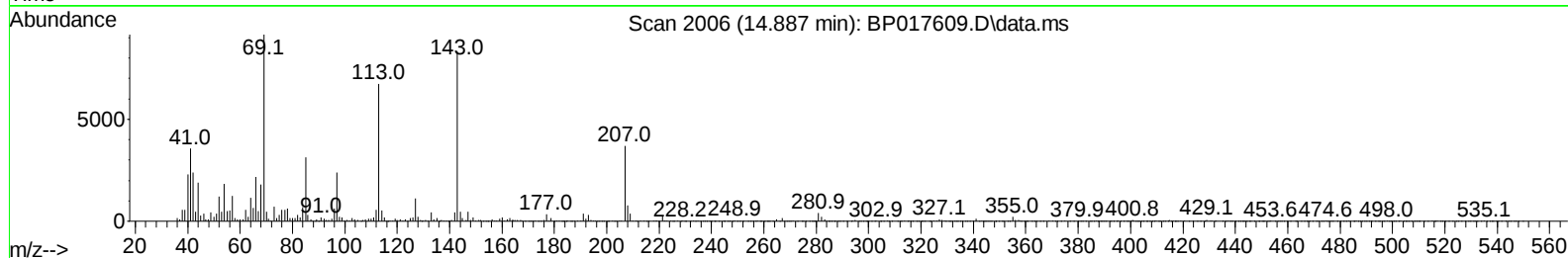
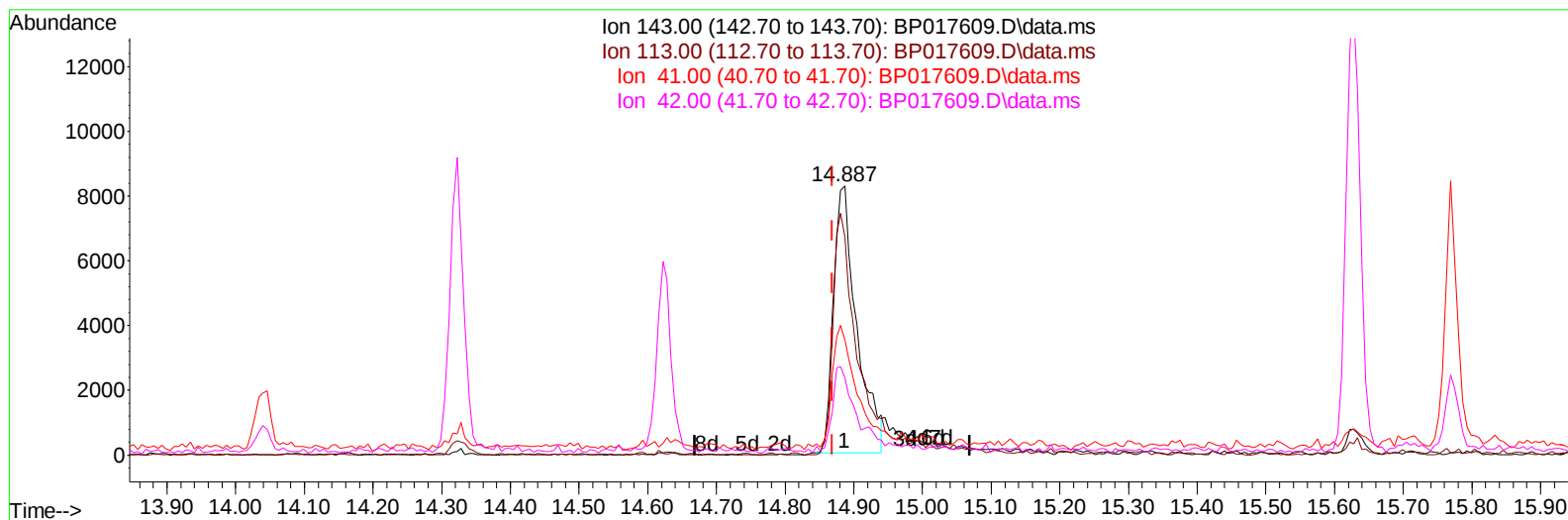
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017609.D
Acq On : 06 Oct 2023 19:21
Operator : MA/JU
Sample : 04624-12
Misc :
ALS Vial : 89 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBJS0

Manual IntegrationsAPPROVED

Quant Time: Oct 06 23:15:26 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 06 05:55:05 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/07/2023
Supervised By :mohammad ahmed 10/07/2023



TIC: BP017609.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.018) 6.20 ng/u1

response 19303

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	81.04
41.00	43.80	42.86
42.00	29.00	28.71

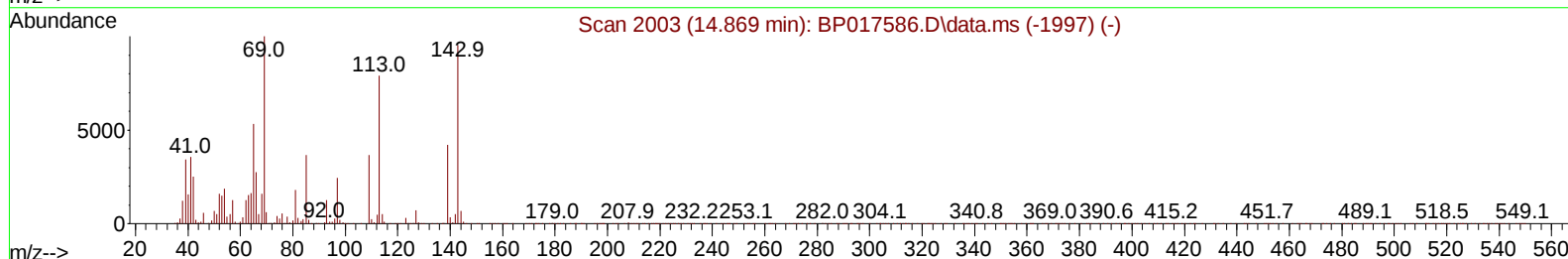
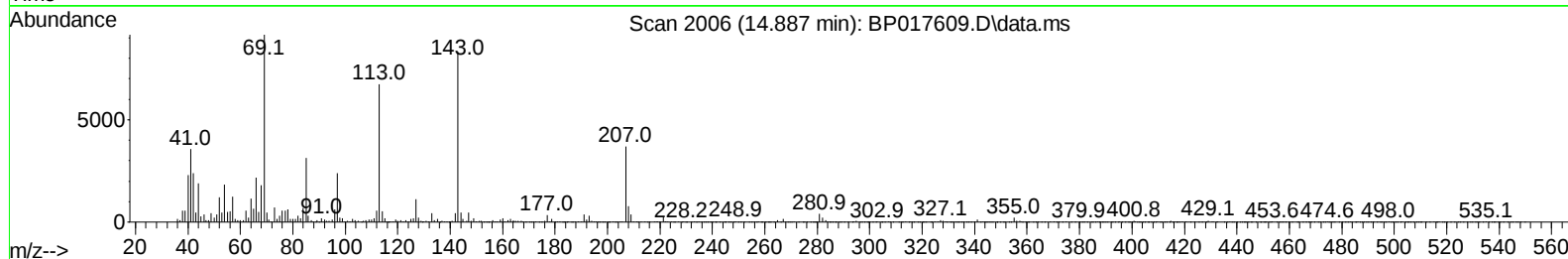
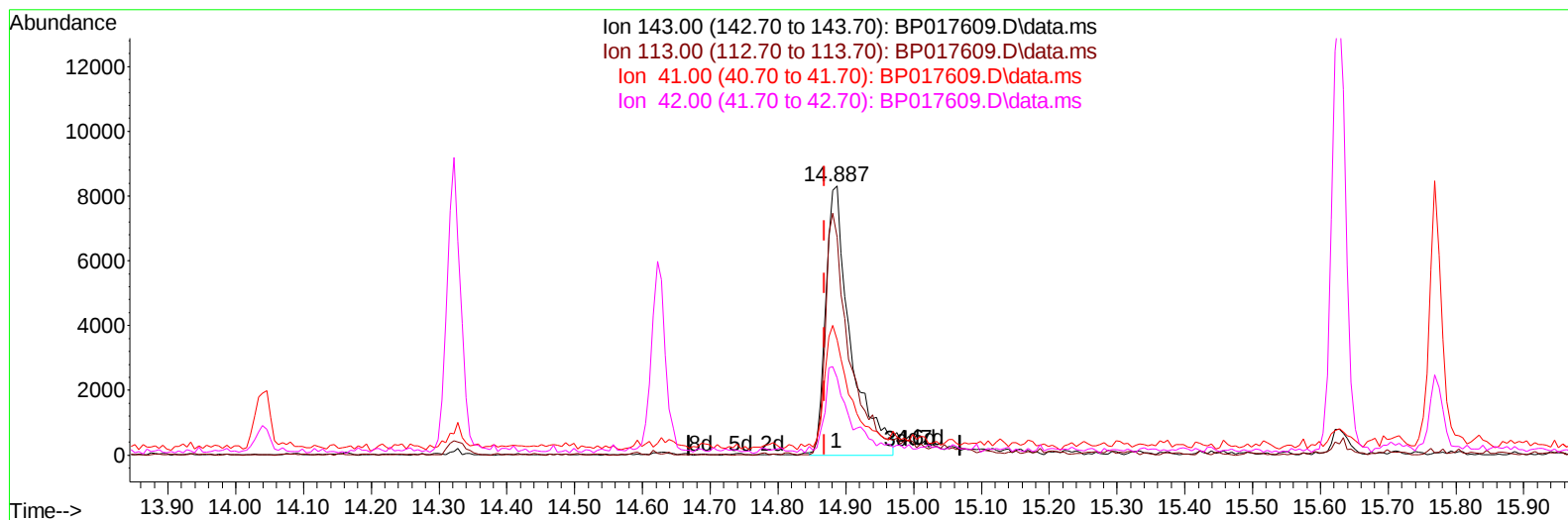
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 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJSO

Manual IntegrationsAPPROVED

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TIC: BP017609.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.018) 6.78 ng/ul m

response 21103

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	81.04
41.00	43.80	42.86
42.00	29.00	28.71

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Instrument :
 BNA_P
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 BBJSO

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/07/2023
 Supervised By :mohammad ahmed 10/07/2023

Quant Time: Oct 06 23:56:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 05:55:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	87918	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	356866	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	228005	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	539322	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	578704	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	651036	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	10793	4.908	ng/uL	0.00
4) Pyridine-d5	3.723	84	82698	14.367	ng/ul	0.00
7) Phenol-d5	7.093	99	67047	9.064	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	185748	41.152	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	190825	32.763	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	124217	20.976	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	115137	41.937	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	123170	39.777	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	218078	37.988	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	266650	32.607	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	827662	46.537	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	864830	44.098	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	21103m	6.776	ng/ul	0.02
60) Fluorene-d10	15.628	176	726657	48.170	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	119075	35.415	ng/ul	0.00
73) Anthracene-d10	17.516	188	1218650	48.815	ng/ul	0.00
81) Pyrene-d10	19.769	212	1575293	48.988	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1635206	49.770	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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